



Full wwPDB X-ray Structure Validation Report

Feb 26, 2014 – 06:18 PM GMT

PDB ID : 2CFH
Title : STRUCTURE OF THE BET3-TPC6B CORE OF TRAPP
Authors : Kummel, D.; Muller, J.J.; Roske, Y.; Henke, N.; Heinemann, U.
Deposited on : 2006-02-21
Resolution : 2.30 Å(reported)

This is a full wwPDB validation report for a publicly released PDB entry.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at <http://wwpdb.org/ValidationPDFNotes.html>

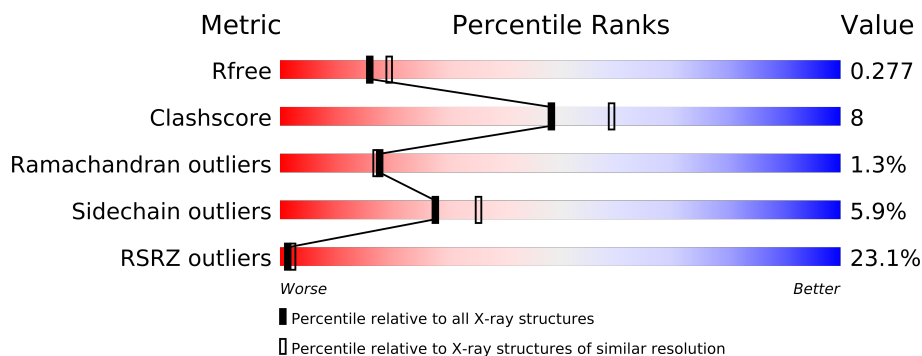
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : dev-1323
EDS : stable22639
Percentile statistics : 21963
Refmac : 5.8.0049
CCP4 : 6.3.0 (Settle)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : stable22683

1 Overall quality at a glance

The reported resolution of this entry is 2.30 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	66092	2929 (2.30-2.30)
Clashscore	79885	3679 (2.30-2.30)
Ramachandran outliers	78287	3642 (2.30-2.30)
Sidechain outliers	78261	3641 (2.30-2.30)
RSRZ outliers	66119	2930 (2.30-2.30)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	194	
1	B	194	
2	C	158	
2	D	158	

The following table lists non-polymeric compounds that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Geometry	Electron density
3	PLM	B	1068	-	X

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 5049 atoms, of which 0 are hydrogen and 0 are deuterium.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called TRAFFICKING PROTEIN PARTICLE COMPLEX SUB-UNIT 3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	162	Total	C	N	O	S	0	0	1
			1276	807	215	246	8			
1	B	162	Total	C	N	O	S	0	0	1
			1276	807	215	246	8			

- Molecule 2 is a protein called TRAFFICKING PROTEIN PARTICLE COMPLEX SUB-UNIT 6B.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	151	Total	C	N	O	S	0	0	1
			1199	764	206	220	9			
2	D	152	Total	C	N	O	S	0	0	1
			1207	769	207	221	10			

- Molecule 3 is PALMITIC ACID (three-letter code: PLM) (formula: C₁₆H₃₂O₂).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			17	16	1		
3	B	1	Total	C	O	0	0
			17	16	1		

- Molecule 4 is water.

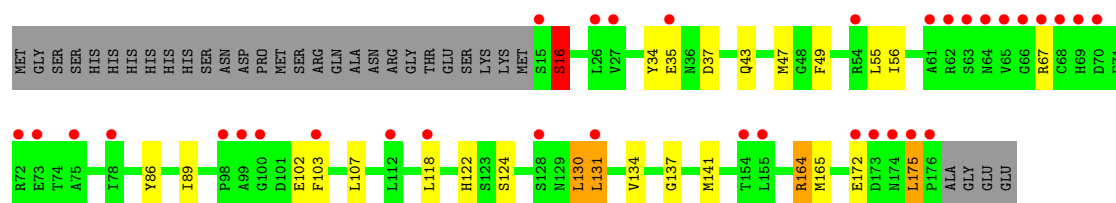
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	18	Total	O	0	0
			18	18		
4	B	12	Total	O	0	0
			12	12		
4	C	10	Total	O	0	0
			10	10		
4	D	17	Total	O	0	0
			17	17		

3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

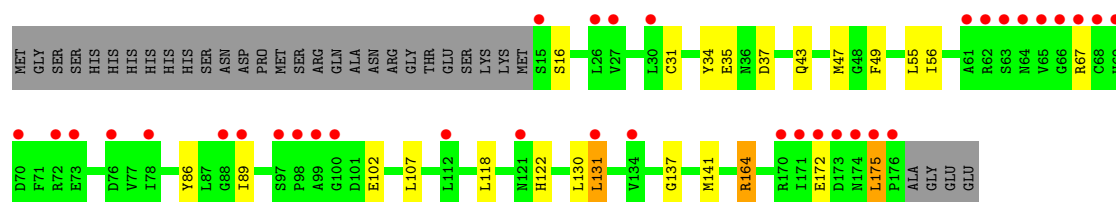
• Molecule 1: TRAFFICKING PROTEIN PARTICLE COMPLEX SUBUNIT 3

Chain A: 



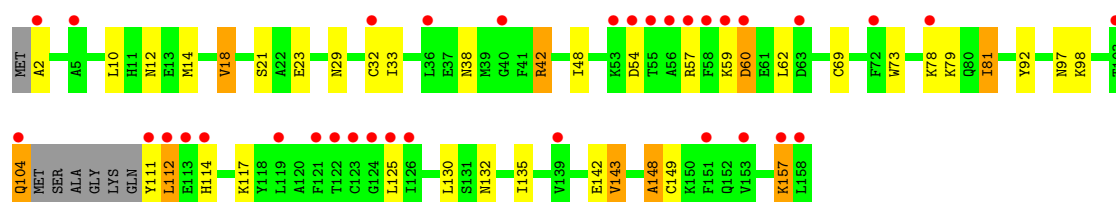
• Molecule 1: TRAFFICKING PROTEIN PARTICLE COMPLEX SUBUNIT 3

Chain B: 



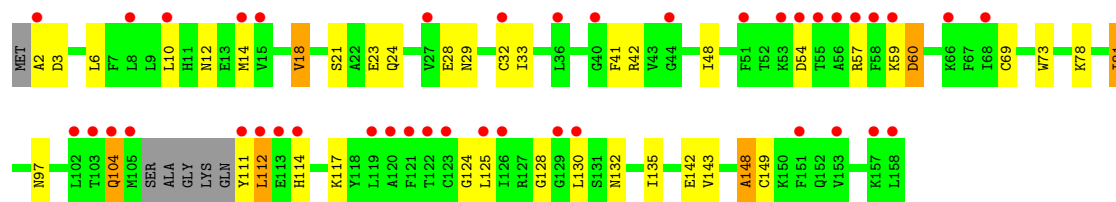
• Molecule 2: TRAFFICKING PROTEIN PARTICLE COMPLEX SUBUNIT 6B

Chain C: 



• Molecule 2: TRAFFICKING PROTEIN PARTICLE COMPLEX SUBUNIT 6B

Chain D: 



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	69.45Å 69.59Å 144.01Å 90.00° 91.55° 90.00°	Depositor
Resolution (Å)	20.00 – 2.30 19.93 – 2.30	Depositor EDS
% Data completeness (in resolution range)	100.0 (20.00-2.30) 98.7 (19.93-2.30)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.85 (at 2.30Å)	Xtriage
Refinement program	REFMAC 5.2.0005	Depositor
R, R_{free}	0.221 , 0.269 0.231 , 0.277	Depositor DCC
R_{free} test set	1513 reflections (5.26%)	DCC
Wilson B-factor (Å ²)	46.6	Xtriage
Anisotropy	0.121	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 47.9	EDS
Estimated twinning fraction	0.000 for k,h,-l 0.000 for -k,-h,-l 0.035 for -h,-k,l	Xtriage
L-test for twinning	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 30262 reflections	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	5049	wwPDB-VP
Average B, all atoms (Å ²)	58.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 34.02 % of the origin peak, indicating pseudo translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo translational symmetry is equal to 7.2959e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹Intensities estimated from amplitudes.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: CSO, PLM

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	0.70	1/1296 (0.1%)	0.72	1/1752 (0.1%)
1	B	0.70	2/1296 (0.2%)	0.72	1/1752 (0.1%)
2	C	0.98	3/1209 (0.2%)	0.87	2/1619 (0.1%)
2	D	0.79	0/1217	0.80	0/1629
All	All	0.80	6/5018 (0.1%)	0.78	4/6752 (0.1%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1
1	B	0	1
2	C	0	1
2	D	0	1
All	All	0	4

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	157	LYS	CE-NZ	20.81	2.01	1.49
2	C	157	LYS	CD-CE	8.59	1.72	1.51
2	C	157	LYS	CG-CD	6.87	1.75	1.52
1	B	175	LEU	C-N	-5.48	1.23	1.34
1	A	175	LEU	C-N	-5.40	1.24	1.34
1	B	31	CYS	CB-SG	-5.07	1.73	1.81

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	157	LYS	CD-CE-NZ	-15.16	76.83	111.70
1	B	67	ARG	C-N-CA	7.19	139.68	121.70
1	A	67	ARG	C-N-CA	7.09	139.43	121.70
2	C	157	LYS	CG-CD-CE	-5.18	96.37	111.90

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	16	SER	Peptide
1	B	16	SER	Peptide
2	C	148	ALA	Peptide
2	D	148	ALA	Peptide

5.2 Close contacts ⓘ

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1276	0	1263	18	0
1	B	1276	0	1263	13	0
2	C	1199	0	1208	30	0
2	D	1207	0	1217	23	0
3	A	17	0	31	1	0
3	B	17	0	31	2	0
4	A	18	0	0	0	0
4	B	12	0	0	0	0
4	C	10	0	0	0	0
4	D	17	0	0	0	0
All	All	5049	0	5013	78	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 8.

All (78) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:C:157:LYS:CD	2:C:157:LYS:CG	1.75	1.57

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:C:32:CSO:SG	2:C:32:CSO:OD	1.92	1.28
2:D:32:CSO:OD	2:D:32:CSO:SG	1.93	1.25
2:C:157:LYS:NZ	2:C:157:LYS:CE	2.01	1.24
2:C:157:LYS:CD	2:C:157:LYS:NZ	2.33	0.92
2:C:69:CYS:HA	2:C:81:ILE:HD11	1.52	0.91
2:D:69:CYS:HA	2:D:81:ILE:HD11	1.51	0.91
2:C:157:LYS:CG	2:C:157:LYS:CE	2.59	0.80
2:C:157:LYS:CD	2:C:157:LYS:CB	2.60	0.79
2:D:21:SER:OG	2:D:29:ASN:ND2	2.27	0.67
2:D:69:CYS:HA	2:D:81:ILE:CD1	2.25	0.65
2:C:21:SER:OG	2:C:29:ASN:ND2	2.31	0.64
2:C:69:CYS:HA	2:C:81:ILE:CD1	2.28	0.62
1:A:16:SER:HB3	2:C:2:ALA:HB3	1.82	0.61
1:B:34:TYR:O	1:B:35:GLU:HG2	2.01	0.61
1:B:55:LEU:HD11	1:B:86:TYR:CE1	2.39	0.57
1:A:89:ILE:HD12	1:A:107:LEU:HD22	1.88	0.56
1:B:55:LEU:HD11	1:B:86:TYR:CZ	2.42	0.55
1:A:56:ILE:HG21	1:A:141:MET:HB2	1.87	0.54
1:A:55:LEU:HD11	1:A:86:TYR:CE1	2.41	0.54
1:B:56:ILE:HG21	1:B:141:MET:HB2	1.89	0.54
2:C:143:VAL:HG23	2:C:148:ALA:O	2.08	0.54
2:D:97:ASN:HA	2:D:148:ALA:HB2	1.90	0.53
1:B:86:TYR:HA	2:D:3:ASP:HB2	1.91	0.51
1:A:16:SER:HB2	2:C:2:ALA:O	2.09	0.51
2:D:69:CYS:SG	2:D:81:ILE:HG13	2.50	0.51
2:C:97:ASN:HA	2:C:148:ALA:HB2	1.92	0.51
1:B:43:GLN:HE21	1:B:47:MET:HG3	1.77	0.50
1:B:89:ILE:HD12	1:B:107:LEU:HD22	1.93	0.50
1:A:34:TYR:O	1:A:35:GLU:HG2	2.12	0.50
1:A:102:GLU:OE1	1:A:164:ARG:NH1	2.45	0.50
2:C:130:LEU:HB3	2:C:135:ILE:HB	1.93	0.50
1:B:37:ASP:OD2	1:B:122:HIS:HA	2.12	0.49
1:A:55:LEU:HD11	1:A:86:TYR:CZ	2.49	0.47
1:B:49:PHE:CE1	1:B:137:GLY:HA2	2.48	0.47
1:A:49:PHE:CE1	1:A:137:GLY:HA2	2.50	0.47
2:C:73:TRP:CG	2:C:81:ILE:HD13	2.49	0.47
1:B:102:GLU:OE1	1:B:164:ARG:NH1	2.48	0.46
2:C:62:LEU:HD22	2:C:92:TYR:OH	2.15	0.46
2:D:143:VAL:HG23	2:D:148:ALA:O	2.15	0.46
2:C:69:CYS:SG	2:C:81:ILE:HG13	2.56	0.46
2:C:79:LYS:NZ	2:C:98:LYS:O	2.43	0.46
2:D:14:MET:HA	2:D:18:VAL:HG23	1.98	0.46

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:D:111:TYR:CD2	2:D:112:LEU:N	2.84	0.46
2:D:130:LEU:HB3	2:D:135:ILE:HB	1.97	0.46
2:D:73:TRP:CG	2:D:81:ILE:HD13	2.52	0.45
1:A:37:ASP:OD2	1:A:122:HIS:HA	2.16	0.45
2:C:111:TYR:CD2	2:C:112:LEU:N	2.85	0.45
2:D:48:ILE:HG21	2:D:132:ASN:HB3	1.99	0.45
2:C:48:ILE:HG21	2:C:132:ASN:HB3	1.98	0.44
2:D:24:GLN:HB2	2:D:28:GLU:OE1	2.18	0.44
1:A:47:MET:CE	2:C:18:VAL:CG2	2.95	0.44
2:C:114:HIS:HB3	2:C:117:LYS:HD2	1.99	0.44
1:A:16:SER:CB	2:C:2:ALA:HB3	2.48	0.44
2:C:38:ASN:O	2:C:42:ARG:NE	2.51	0.43
1:A:131:LEU:HD21	3:A:1068:PLM:HC2	1.98	0.43
2:D:12:ASN:HB2	2:D:104:GLN:HG3	2.00	0.43
2:C:60:ASP:OD1	2:C:60:ASP:C	2.56	0.43
2:D:33:ILE:HD13	2:D:33:ILE:HA	1.85	0.43
1:A:103:PHE:CZ	1:A:165:MET:HG3	2.53	0.43
2:C:59:LYS:O	2:C:60:ASP:HB3	2.19	0.43
2:C:33:ILE:HA	2:C:33:ILE:HD13	1.86	0.43
2:C:14:MET:HA	2:C:18:VAL:HG23	2.01	0.42
2:C:12:ASN:HB2	2:C:104:GLN:HG3	2.00	0.42
2:D:41:PHE:CE1	2:D:128:GLY:HA2	2.55	0.42
2:C:23:GLU:HA	2:C:29:ASN:HD21	1.85	0.42
2:D:23:GLU:HA	2:D:29:ASN:HD21	1.84	0.41
1:B:86:TYR:O	2:D:2:ALA:HA	2.21	0.41
1:A:130:LEU:O	1:A:134:VAL:HG23	2.20	0.41
2:D:41:PHE:HA	2:D:124:GLY:O	2.20	0.41
1:B:131:LEU:HD21	3:B:1068:PLM:HC2	2.02	0.41
3:B:1068:PLM:HF1	2:D:6:LEU:HD11	2.02	0.41
2:D:59:LYS:O	2:D:60:ASP:HB3	2.21	0.41
1:A:37:ASP:HB2	1:A:124:SER:HB2	2.02	0.41
1:A:43:GLN:NE2	1:A:43:GLN:HA	2.35	0.40
1:A:164:ARG:HA	1:A:164:ARG:HD3	1.84	0.40
2:D:114:HIS:HB3	2:D:117:LYS:HD2	2.03	0.40
1:B:164:ARG:HA	1:B:164:ARG:HD3	1.85	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone ⓘ

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution. The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	160/194 (82%)	154 (96%)	5 (3%)	1 (1%)	33	39
1	B	160/194 (82%)	154 (96%)	5 (3%)	1 (1%)	33	39
2	C	146/158 (92%)	132 (90%)	11 (8%)	3 (2%)	11	8
2	D	147/158 (93%)	131 (89%)	13 (9%)	3 (2%)	11	8
All	All	613/704 (87%)	571 (93%)	34 (6%)	8 (1%)	18	17

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	C	60	ASP
2	C	149	CYS
2	D	60	ASP
2	D	149	CYS
2	C	112	LEU
2	D	112	LEU
1	A	175	LEU
1	B	175	LEU

5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution. The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	142/170 (84%)	136 (96%)	6 (4%)	40	53
1	B	142/170 (84%)	137 (96%)	5 (4%)	48	63
2	C	129/136 (95%)	118 (92%)	11 (8%)	15	18
2	D	130/136 (96%)	120 (92%)	10 (8%)	18	22
All	All	543/612 (89%)	511 (94%)	32 (6%)	28	35

All (32) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	16	SER
1	A	118	LEU
1	A	130	LEU
1	A	131	LEU
1	A	164	ARG
1	A	172	GLU
1	B	118	LEU
1	B	130	LEU
1	B	131	LEU
1	B	164	ARG
1	B	172	GLU
2	C	10	LEU
2	C	18	VAL
2	C	42	ARG
2	C	54	ASP
2	C	57	ARG
2	C	78	LYS
2	C	81	ILE
2	C	104	GLN
2	C	125	LEU
2	C	142	GLU
2	C	143	VAL
2	D	10	LEU
2	D	18	VAL
2	D	42	ARG
2	D	54	ASP
2	D	57	ARG
2	D	78	LYS
2	D	81	ILE
2	D	104	GLN
2	D	125	LEU
2	D	142	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	43	GLN
1	B	43	GLN
2	C	29	ASN
2	C	38	ASN
2	C	83	ASN

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Mol	Chain	Res	Type
2	C	95	GLN
2	D	29	ASN
2	D	38	ASN
2	D	83	ASN
2	D	95	GLN

5.3.3 RNA ⓘ

There are no RNA chains in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

2 non-standard protein/DNA/RNA residues are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	CSO	C	32	2	6,6,7	7.95	4 (66%)	3,6,8	2.25	2 (66%)
2	CSO	D	32	2	6,6,7	7.36	4 (66%)	3,6,8	2.03	1 (33%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the chemical component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	CSO	C	32	2	-	0/2/5/7	0/0/0/0
2	CSO	D	32	2	-	0/2/5/7	0/0/0/0

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	32	CSO	O-C	16.81	1.23	1.11
2	D	32	CSO	O-C	15.18	1.21	1.11

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	32	CSO	OD-SG	7.27	1.93	1.62
2	C	32	CSO	OD-SG	6.97	1.92	1.62
2	C	32	CSO	CB-SG	-6.35	1.74	1.82
2	D	32	CSO	CB-SG	-5.98	1.75	1.82
2	C	32	CSO	CA-C	2.45	1.52	1.48
2	D	32	CSO	CA-C	2.19	1.52	1.48

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	32	CSO	CA-CB-SG	-2.57	108.45	113.01
2	D	32	CSO	CA-CB-SG	-2.38	108.78	113.01
2	C	32	CSO	CB-CA-N	-2.24	106.47	110.27

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

2 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
3	PLM	A	1068	1	16,16,17	4.54	2 (12%)	13,15,17	0.57	0
3	PLM	B	1068	1	16,16,17	4.80	2 (12%)	13,15,17	0.61	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral

centers analysed, the number of these observed in the model and the number defined in the chemical component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	PLM	A	1068	1	-	0/13/14/15	0/0/0/0
3	PLM	B	1068	1	-	0/13/14/15	0/0/0/0

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	1068	PLM	O2-C1	19.00	1.24	1.11
3	A	1068	PLM	O2-C1	17.96	1.23	1.11
3	B	1068	PLM	C2-C1	2.26	1.54	1.47
3	A	1068	PLM	C2-C1	2.08	1.53	1.47

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	162/194 (83%)	1.31	34 (20%) 1 3	57, 59, 60, 61	0
1	B	162/194 (83%)	1.40	35 (21%) 1 2	57, 59, 61, 62	0
2	C	151/158 (95%)	1.29	34 (22%) 1 2	54, 59, 61, 64	0
2	D	152/158 (96%)	1.42	41 (26%) 1 1	54, 59, 61, 65	0
All	All	627/704 (89%)	1.35	144 (22%) 1 2	54, 59, 61, 65	0

All (144) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	176	PRO	10.5
1	B	176	PRO	10.0
1	A	175	LEU	10.0
1	B	175	LEU	8.5
2	C	111	TYR	7.5
2	D	56	ALA	7.2
2	D	111	TYR	6.2
2	D	57	ARG	5.4
2	D	105	MET	5.4
1	B	65	VAL	5.4
1	B	67	ARG	5.3
2	C	112	LEU	5.0
1	A	99	ALA	5.0
1	A	69	HIS	5.0
1	A	98	PRO	5.0
2	C	103	THR	4.8
1	B	69	HIS	4.7
1	B	174	ASN	4.7
1	B	172	GLU	4.5
2	D	158	LEU	4.5
1	B	68	CYS	4.5

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Mol	Chain	Res	Type	RSRZ
1	A	15	SER	4.4
2	C	55	THR	4.4
2	D	125	LEU	4.3
2	C	57	ARG	4.2
1	B	66	GLY	4.2
1	B	73	GLU	4.1
1	B	70	ASP	4.1
2	D	123	CYS	4.1
2	D	58	PHE	4.1
2	C	56	ALA	4.0
1	A	100	GLY	4.0
2	D	103	THR	3.9
2	C	125	LEU	3.9
2	C	53	LYS	3.9
1	A	65	VAL	3.9
1	A	27	VAL	3.8
2	C	126	ILE	3.8
1	A	174	ASN	3.8
2	C	59	LYS	3.7
2	C	122	THR	3.6
2	D	157	LYS	3.6
1	A	67	ARG	3.6
2	D	126	ILE	3.6
1	A	70	ASP	3.5
1	A	61	ALA	3.5
1	A	73	GLU	3.5
2	D	113	GLU	3.5
2	D	53	LYS	3.5
1	B	173	ASP	3.4
1	B	64	ASN	3.4
1	A	54	ARG	3.3
1	A	63	SER	3.3
1	B	27	VAL	3.3
1	B	170	ARG	3.3
2	C	58	PHE	3.2
1	B	99	ALA	3.2
1	A	64	ASN	3.2
2	D	36	LEU	3.1
2	D	55	THR	3.1
2	D	122	THR	3.1
1	B	100	GLY	3.1
2	D	104	GLN	3.0

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Mol	Chain	Res	Type	RSRZ
2	D	120	ALA	3.0
2	D	130	LEU	3.0
2	C	157	LYS	3.0
1	B	98	PRO	2.9
1	A	68	CYS	2.9
2	D	2	ALA	2.9
1	B	26	LEU	2.8
2	C	36	LEU	2.8
2	C	113	GLU	2.8
1	A	103	PHE	2.8
1	A	26	LEU	2.8
2	D	151	PHE	2.8
1	A	118	LEU	2.8
2	C	72	PHE	2.7
2	D	119	LEU	2.7
2	C	54	ASP	2.7
2	D	10	LEU	2.7
2	D	54	ASP	2.7
2	C	121	PHE	2.7
1	B	88	GLY	2.7
1	A	75	ALA	2.7
1	A	131	LEU	2.6
2	C	123	CYS	2.6
1	B	89	ILE	2.6
2	C	2	ALA	2.6
1	B	131	LEU	2.6
1	A	128	SER	2.6
1	B	72	ARG	2.6
1	A	172	GLU	2.5
1	A	173	ASP	2.5
1	A	78	ILE	2.5
2	D	66	LYS	2.5
1	B	97	SER	2.5
2	D	68	ILE	2.5
2	C	158	LEU	2.4
1	B	76	ASP	2.4
2	D	51	PHE	2.4
2	D	14	MET	2.4
2	D	112	LEU	2.4
2	C	124	GLY	2.4
1	B	121	ASN	2.4
2	C	104	GLN	2.4

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Mol	Chain	Res	Type	RSRZ
2	D	8	LEU	2.4
2	D	114	HIS	2.3
2	D	44	GLY	2.3
2	D	15	VAL	2.3
2	D	129	GLY	2.3
2	D	59	LYS	2.3
2	C	78	LYS	2.3
2	D	40	GLY	2.3
2	D	32	CSO	2.3
2	C	151	PHE	2.2
1	A	62	ARG	2.2
2	C	63	ASP	2.2
1	A	66	GLY	2.2
1	B	30	LEU	2.2
2	C	40	GLY	2.2
2	D	121	PHE	2.2
2	D	153	VAL	2.2
2	C	32	CSO	2.1
1	A	155	LEU	2.1
2	C	139	VAL	2.1
1	A	35	GLU	2.1
2	C	60	ASP	2.1
1	B	134	VAL	2.1
1	A	72	ARG	2.1
1	B	15	SER	2.1
2	C	5	ALA	2.1
1	B	112	LEU	2.1
1	B	63	SER	2.1
1	B	61	ALA	2.1
2	C	119	LEU	2.1
1	B	78	ILE	2.1
2	D	102	LEU	2.0
2	C	153	VAL	2.0
1	B	171	ILE	2.0
2	C	114	HIS	2.0
1	A	112	LEU	2.0
1	B	62	ARG	2.0
2	D	27	VAL	2.0
1	A	154	THR	2.0

6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSR	LLDF	B-factors(Å ²)	Q<0.9
2	CSO	C	32	7/8	0.19	0.13	48,57,57,58	0
2	CSO	D	32	7/8	0.20	-0.18	48,57,57,57	0

6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

6.4 Ligands ⓘ

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSR	LLDF	B-factors(Å ²)	Q<0.9
3	PLM	B	1068	17/18	0.47	2.60	58,59,61,61	0
3	PLM	A	1068	17/18	0.27	0.83	58,59,61,61	0

6.5 Other polymers ⓘ

There are no such residues in this entry.